```
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LOGINID:SSPTANXR1625

PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2


* * * * * * * * * *  Welcome to STN International  * * * * * * * * * *

 NEWS  1          Web Page for STN Seminar Schedule - N. America
 NEWS  2  NOV 21  CAS patent coverage to include exemplified prophetic
                  substances identified in English-, French-, German-,
                  and Japanese-language basic patents from 2004-present
 NEWS  3  NOV 26  MARPAT enhanced with FSORT command
 NEWS  4  NOV 26  CHEMSAFE now available on STN Easy
 NEWS  5  NOV 26  Two new SET commands increase convenience of STN
                  searching
 NEWS  6  DEC 01  ChemPort single article sales feature unavailable
 NEWS  7  DEC 12  GBFULL now offers single source for full-text
                  coverage of complete UK patent families
 NEWS  8  DEC 17  Fifty-one pharmaceutical ingredients added to PS
 NEWS  9  JAN 06  The retention policy for unread STNmail messages
                  will change in 2009 for STN-Columbus and STN-Tokyo
 NEWS 10  JAN 07  WPIDS, WPINDEX, and WPIX enhanced Japanese Patent
                  Classification Data
 NEWS 11  FEB 02  Simultaneous left and right truncation (SLART) added
                  for CERAB, COMPUAB, ELCOM, and SOLIDSTATEM
 NEWS 12  FEB 02  GENBANK enhanced with SET PLURALS and SET SPELLING

 NEWS EXPRESS JUNE 27 08 CURRENT WINDOWS VERSION IS V8.3,
              AND CURRENT DISCOVER FILE IS DATED 23 JUNE 2008.

 NEWS HOURS    STN Operating Hours Plus Help Desk Availability
 NEWS LOGIN    Welcome Banner and News Items
 NEWS IPC8     For general information regarding STN implementation of IPC 8

Enter NEWS followed by the item number or name to see news on that
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* * * * * * * * * * * * * * * * STN  Columbus  * * * * * * * * * * * * * * * *

FILE 'HOME' ENTERED AT 14:55:22 ON 04 FEB 2009

=> file reg
COST IN U.S. DOLLARS                                SINCE FILE      TOTAL
                                                       ENTRY    SESSION
FULL ESTIMATED COST                                     0.22       0.22

FILE 'REGISTRY' ENTERED AT 14:55:31 ON 04 FEB 2009
```

Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.

STRUCTURE FILE UPDATES: 3 FEB 2009 HIGHEST RN 1100396-01-7
DICTIONARY FILE UPDATES: 3 FEB 2009 HIGHEST RN 1100396-01-7

New CAS Information Use Policies, enter HELP USAGETERMS for details.

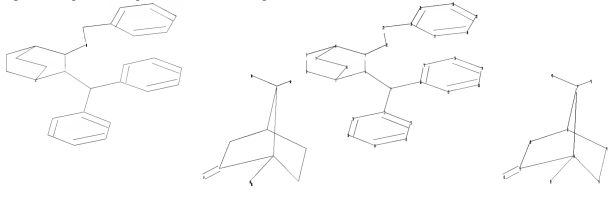
TSCA INFORMATION NOW CURRENT THROUGH July 5, 2008.

  Please note that search-term pricing does apply when
  conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:

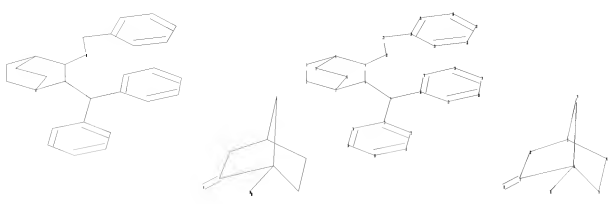
http://www.cas.org/support/stngen/stndoc/properties.html

=>
Uploading C:\Program Files\Stnexp\Queries\10588049.str



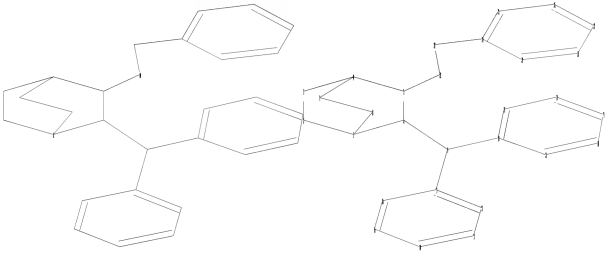
chain nodes :
9 22 23 37 38 40 41

```
ring nodes :
1  2  3  4  5  6  7  8  10  11  12  13  14  15  16  17  18  19  20  21  24  25  26
27  28  29  31  32  33  34  35  36  39
chain bonds :
5-22  6-9  9-10  9-11  22-23  23-24  31-37  36-38  39-40  39-41
ring bonds :
1-2  1-6  1-8  2-3  3-4  4-5  4-7  5-6  7-8  10-17  10-21  11-12  11-16  12-13
13-14  14-15  15-16  17-18  18-19  19-20  20-21  24-25  24-29  25-26  26-27  27-28
28-29  31-32  31-36  32-33  33-34  33-39  34-35  35-36  36-39
exact/norm bonds :
1-2  1-6  1-8  2-3  3-4  4-5  4-7  5-6  5-22  7-8  22-23  31-37
exact bonds :
6-9  9-10  9-11  23-24  31-32  31-36  32-33  33-34  33-39  34-35  35-36  36-38
36-39  39-40  39-41
normalized bonds :
10-17  10-21  11-12  11-16  12-13  13-14  14-15  15-16  17-18  18-19  19-20  20-21
24-25  24-29  25-26  26-27  27-28  28-29
isolated ring systems :
containing 1 : 10 : 11 : 24 : 31 :


Match level :
1:Atom  2:Atom  3:Atom  4:Atom  5:Atom  6:Atom  7:Atom  8:Atom  9:CLASS  10:Atom
11:Atom  12:Atom  13:Atom  14:Atom  15:Atom  16:Atom  17:Atom  18:Atom  19:Atom
20:Atom  21:Atom  22:CLASS  23:CLASS  24:Atom  25:Atom  26:Atom  27:Atom  28:Atom
29:Atom  31:Atom  32:Atom  33:Atom  34:Atom  35:Atom  36:Atom  37:CLASS  38:CLASS
39:Atom  40:CLASS  41:CLASS



L1      STRUCTURE UPLOADED

=> d l1
L1 HAS NO ANSWERS
L1              STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY -  AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.


=> s l1 full
FULL SEARCH INITIATED 14:55:54 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -        0 TO ITERATE

100.0% PROCESSED        0 ITERATIONS                                 0 ANSWERS
SEARCH TIME: 00.00.01

L2            0 SEA SSS FUL L1

=>
Uploading C:\Program Files\Stnexp\Queries\10588049a.str
```



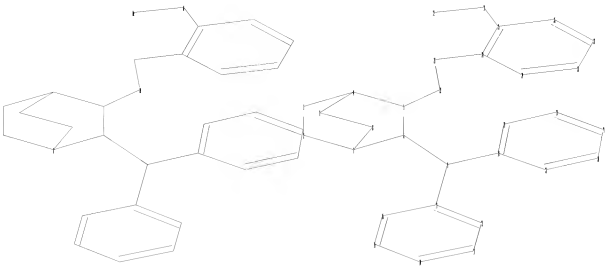
```
chain nodes :
9  22  23  37  38
ring nodes :
1  2  3  4  5  6  7  8  10  11  12  13  14  15  16  17  18  19  20  21  24  25  26
27  28  29  31  32  33  34  35  36  39
chain bonds :
5-22  6-9  9-10  9-11  22-23  23-24  31-37  36-38
ring bonds :
1-2  1-6  1-8  2-3  3-4  4-5  4-7  5-6  7-8  10-17  10-21  11-12  11-16  12-13
13-14  14-15  15-16  17-18  18-19  19-20  20-21  24-25  24-29  25-26  26-27  27-28
28-29  31-32  31-36  32-33  33-34  33-39  34-35  35-36  36-39
exact/norm bonds :
1-2  1-6  1-8  2-3  3-4  4-5  4-7  5-6  5-22  7-8  22-23  31-37
exact bonds :
6-9  9-10  9-11  23-24  31-32  31-36  32-33  33-34  33-39  34-35  35-36  36-38
36-39
normalized bonds :
10-17  10-21  11-12  11-16  12-13  13-14  14-15  15-16  17-18  18-19  19-20  20-21
24-25  24-29  25-26  26-27  27-28  28-29
isolated ring systems :
containing 1 : 10 : 11 : 24 : 31 :


Match level :
1:Atom  2:Atom  3:Atom  4:Atom  5:Atom  6:Atom  7:Atom  8:Atom  9:CLASS  10:Atom
11:Atom  12:Atom  13:Atom  14:Atom  15:Atom  16:Atom  17:Atom  18:Atom  19:Atom
20:Atom  21:Atom  22:CLASS  23:CLASS  24:Atom  25:Atom  26:Atom  27:Atom  28:Atom
29:Atom  31:Atom  32:Atom  33:Atom  34:Atom  35:Atom  36:Atom  37:CLASS  38:CLASS
39:Atom
```

```
L3      STRUCTURE UPLOADED

=> d l3
L3 HAS NO ANSWERS
L3              STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY -  AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.


=> s l3 full
FULL SEARCH INITIATED 14:57:30 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -        9 TO ITERATE

100.0% PROCESSED        9 ITERATIONS                                    0 ANSWERS
SEARCH TIME: 00.00.01

L4            0 SEA SSS FUL L3

=>
Uploading C:\Program Files\Stnexp\Queries\10588049b.str
```



```
chain nodes :
9  22  23
ring nodes :
1  2  3  4  5  6  7  8  10  11  12  13  14  15  16  17  18  19  20  21  24  25  26
27  28  29
chain bonds :
5-22  6-9  9-10  9-11  22-23  23-24
ring bonds :
1-2  1-6  1-8  2-3  3-4  4-5  4-7  5-6  7-8  10-17  10-21  11-12  11-16  12-13
13-14  14-15  15-16  17-18  18-19  19-20  20-21  24-25  24-29  25-26  26-27  27-28
28-29
```

```
exact/norm bonds :
1-2  1-6  1-8  2-3  3-4  4-5  4-7  5-6  5-22  7-8  22-23
exact bonds :
6-9  9-10  9-11  23-24
normalized bonds :
10-17  10-21  11-12  11-16  12-13  13-14  14-15  15-16  17-18  18-19  19-20  20-21
24-25  24-29  25-26  26-27  27-28  28-29
isolated ring systems :
containing 1 : 10 : 11 : 24 :


Match level :
1:Atom  2:Atom  3:Atom  4:Atom  5:Atom  6:Atom  7:Atom  8:Atom  9:CLASS  10:Atom
11:Atom  12:Atom  13:Atom  14:Atom  15:Atom  16:Atom  17:Atom  18:Atom  19:Atom
20:Atom  21:Atom  22:CLASS  23:CLASS  24:Atom  25:Atom  26:Atom  27:Atom  28:Atom
29:Atom



L5      STRUCTURE UPLOADED

=> d l5
L5 HAS NO ANSWERS
L5              STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY -  AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.


=> s l5 full
FULL SEARCH INITIATED 14:58:31 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -     1044 TO ITERATE

100.0% PROCESSED     1044 ITERATIONS                              431 ANSWERS
SEARCH TIME: 00.00.01

L6          431 SEA SSS FUL L5

=>
Uploading C:\Program Files\Stnexp\Queries\10588049c.str
```



chain nodes :
9 22 23 31 32
ring nodes :
1 2 3 4 5 6 7 8 10 11 12 13 14 15 16 17 18 19 20 21 24 25 26 27 28 29
chain bonds :
5-22 6-9 9-10 9-11 22-23 23-24 25-31 31-32
ring bonds :
1-2 1-6 1-8 2-3 3-4 4-5 4-7 5-6 7-8 10-17 10-21 11-12 11-16 12-13 13-14 14-15 15-16 17-18 18-19 19-20 20-21 24-25 24-29 25-26 26-27 27-28 28-29
exact/norm bonds :
1-2 1-6 1-8 2-3 3-4 4-5 4-7 5-6 5-22 7-8 22-23 25-31
exact bonds :
6-9 9-10 9-11 23-24 31-32
normalized bonds :
10-17 10-21 11-12 11-16 12-13 13-14 14-15 15-16 17-18 18-19 19-20 20-21 24-25 24-29 25-26 26-27 27-28 28-29
isolated ring systems :
containing 1 : 10 : 11 : 24 :

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:CLASS 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom 21:Atom 22:CLASS 23:CLASS 24:Atom 25:Atom 26:Atom 27:Atom 28:Atom 29:Atom 31:CLASS 32:CLASS

L7 STRUCTURE UPLOADED

```
=> d l7
L7 HAS NO ANSWERS
L7              STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY -  AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.


=> s l7 full
FULL SEARCH INITIATED 15:02:48 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -      500 TO ITERATE

100.0% PROCESSED      500 ITERATIONS                              335 ANSWERS
SEARCH TIME: 00.00.01

L8          335 SEA SSS FUL L7

=> file casreact
COST IN U.S. DOLLARS                                SINCE FILE      TOTAL
                                                       ENTRY    SESSION
FULL ESTIMATED COST                                   748.32     748.54

FILE 'CASREACT' ENTERED AT 15:03:47 ON 04 FEB 2009
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FILE CONTENT:1840 - 1 Feb 2009 VOL 150 ISS 6

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*                                                          *
*     CASREACT now has more than 16.5 million reactions    *
*                                                          *
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Biotransformations database compiled under the direction of
Professor Dr. Klaus Kieslich; organic reactions, portions copyright
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This file contains CAS Registry Numbers for easy and accurate substance
identification.

=>
Uploading C:\Program Files\Stnexp\Queries\10588049d.str
```
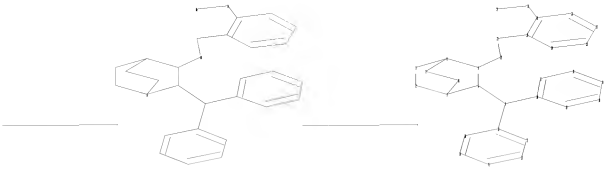



chain nodes :
9 22 23 31 32
ring nodes :
1 2 3 4 5 6 7 8 10 11 12 13 14 15 16 17 18 19 20 21 24 25 26 27 28 29
chain bonds :
5-22 6-9 9-10 9-11 22-23 23-24 25-31 31-32
ring bonds :
1-2 1-6 1-8 2-3 3-4 4-5 4-7 5-6 7-8 10-17 10-21 11-12 11-16 12-13 13-14 14-15 15-16 17-18 18-19 19-20 20-21 24-25 24-29 25-26 26-27 27-28 28-29
exact/norm bonds :
1-2 1-6 1-8 2-3 3-4 4-5 4-7 5-6 5-22 7-8 22-23 25-31
exact bonds :
6-9 9-10 9-11 23-24 31-32
normalized bonds :
10-17 10-21 11-12 11-16 12-13 13-14 14-15 15-16 17-18 18-19 19-20 20-21 24-25 24-29 25-26 26-27 27-28 28-29
isolated ring systems :
containing 1 : 10 : 11 : 24 :

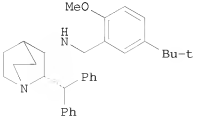
Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:CLASS 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom 21:Atom 22:CLASS 23:CLASS 24:Atom 25:Atom 26:Atom 27:Atom 28:Atom 29:Atom 31:CLASS 32:CLASS
fragments assigned product role:
containing 1

```
L9      STRUCTURE UPLOADED

=> d l9
L9 HAS NO ANSWERS
L9              STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY -  AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.


=> s l9 full
FULL SEARCH INITIATED 15:04:09 FILE 'CASREACT'
SCREENING COMPLETE -       64 REACTIONS TO VERIFY FROM       10 DOCUMENTS

100.0% DONE      64 VERIFIED      35 HIT RXNS                    7 DOCS
SEARCH TIME: 00.00.01

L10           7 SEA SSS FUL L9 (    35 REACTIONS)

=> d ibib abs fhit tot
```

L10 ANSWER 1 OF 7 CASREACT COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 143:230050 CASREACT
TITLE: Process for preparation of 1-(2S,3S)-2-benzhydryl-N-(5-tert-butyl-2-methoxybenzyl)quinuclidin-3-amine
INVENTOR(S): Basford, Patricia Ann; Post, Ronald James; Smith, Julian Duncan; Taber, Geraldine Patricia
PATENT ASSIGNEE(S): Pfizer Products Inc., USA
SOURCE: PCT Int. Appl., 43 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2005075473	A1	20050818	WO 2005-IB221	20050126
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
AU 2005210259	A1	20050818	AU 2005-210259	20050126
CA 2554360	A1	20050818	CA 2005-2554360	20050126
EP 1713801	A1	20061025	EP 2005-702373	20050126
EP 1713801	B1	20071212		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, FI, RO, CY, TR, BG, CZ, EE, HU, PL, SK, IS				
CN 1914202	A	20070214	CN 2005-80003898	20050126
BR 2005007334	A	20070703	BR 2005-7334	20050126
JP 2007519710	T	20070719	JP 2006-550351	20050126
AT 380811	T	20071215	AT 2005-702373	20050126
RU 2320659	C1	20080327	RU 2006-127969	20050126
ES 2296131	T3	20080416	ES 2005-702373	20050126
NO 2006003268	A	20060815	NO 2006-3268	20060713
IN 2006DN04065	A	20070713	IN 2006-DN4065	20060714
MX 2006008441	A	20061002	MX 2006-8441	20060726
KR 2006127966	A	20061213	KR 2006-715573	20060801
KR 812046	B1	20080310		
PRIORITY APPLN. INFO.:			US 2004-541323P	20040202
			WO 2005-IB221	20050126

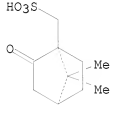
OTHER SOURCE(S): MARPAT 143:230050
GI



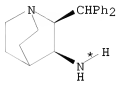
I

AB This invention relates to an improved process for the preparation and purification of (2S,3S)-2-benzhydryl-N-(5-tert-butyl-2-methoxybenzyl)quinuclidin-3-amine (I), which is useful as an antiemetic agent (no biol. testing data), and its pharmaceutically acceptable salts. In particular, the invention is directed to an improved synthesis of the monohydrate monocitrate salt of I.

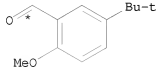
RX(1) OF 6 A + B ===> C...



A: CM 1

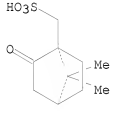


A: CM 2

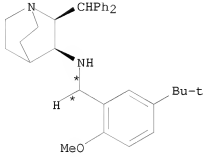


B

(1)



C: CM 1



C: CM 2

RX(1) RCT A 862543-53-1

STAGE(1)
RGT D 1333-74-0 H2
CAT 7440-05-3 Pd
SOL 7732-18-5 Water, 67-63-0 Me2CHOH
CON 4 hours, 75 - 80 deg C, 50 psi

STAGE(2)
RCT B 85943-26-6
SOL 67-63-0 Me2CHOH
CON 2 hours, 75 - 80 deg C

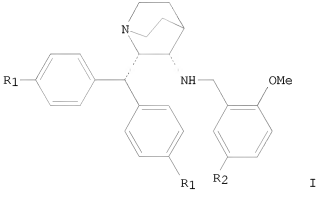
STAGE(3)
RGT D 1333-74-0 H2
CON SUBSTAGE(1) 30 - 40 deg C
SUBSTAGE(2) 3.5 hours, 75 - 80 deg C, 50 psi
SUBSTAGE(3) 10 hours, 25 - 30 deg C, 10 psi
SUBSTAGE(4) 11.5 hours, 75 - 80 deg C, 50 psi
SUBSTAGE(5) 10 hours, 25 - 30 deg C, 10 psi

SUBSTAGE(6) 3 hours, 75 - 80 deg C, 50 psi

PRO C 862543-52-0

REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

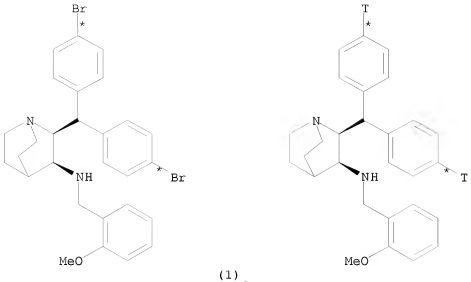
L10 ANSWER 2 OF 7 CASREACT COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 141:6864 CASREACT
TITLE: Tritiation of nonpeptide substance P antagonist CP-96,345 and its azido analogue. Synthetic and characterization details
AUTHOR(S): Egan, Judith A.; Filer, Crist N.
CORPORATE SOURCE: PerkinElmer Life and Analytical Sciences, Inc., Boston, MA, 02118, USA
SOURCE: Applied Radiation and Isotopes (2003), 59(5-6), 333-335
CODEN: ARISEF; ISSN: 0969-8043
PUBLISHER: Elsevier Science B.V.
DOCUMENT TYPE: Journal
LANGUAGE: English
GI



I

AB CP-96,345 was the first nonpeptide antagonist discovered for the SP receptor and [3H] CP-96,345 was required to study the mechanism of receptor action. The radioligand I (R1 = T, R2 = H) was prepared at high specific activity by catalytic dehalogenation of a dibrominated precursor I (R1 = Br, R2 = H). The photoaffinity analog I (R1 = T, R2 = N3) was also prepared from precursor I (R1 = Br, R2 = NH2) using the same approach followed by diazotization and azidation with NaN3.

RX(1) OF 4 A ===> B



A B

RX(1) RCT A 135007-76-0
RGT C 10028-17-8 Tritium
PRO B 135007-77-1
CAT 7440-05-3 Pd
SOL 64-17-5 EtOH
CON 1 hour, room temperature

REFERENCE COUNT: 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L10 ANSWER 3 OF 7 CASREACT COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 137:369932 CASREACT

TITLE: Cooperative problem solving: investigation into the oxidative degradation of CJ-11,974-01 and [14C]CJ-11,974-01

AUTHOR(S): Zandi, Kathleen S.; Huff, Barbara B.; Kamel, Amin; Larmann, John; Massefski, Walter W.; McCarthy, Keith E.; Miller, Sandra A.; Smith, Scott W.

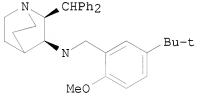
CORPORATE SOURCE: Radiochemical Synthesis, Pfizer Central Research, Groton, CT, 06340, USA

SOURCE: Synthesis and Applications of Isotopically Labelled Compounds, Proceedings of the International Symposium, 7th, Dresden, Germany, June 18-22, 2000 (2001), Meeting Date 2000, 232-235. Editor(s): Pleiss, Ulrich; Voges, Rolf. John Wiley & Sons Ltd.: Chichester, UK.
CODEN: 69CIJC; ISBN: 0-471-49501-8

DOCUMENT TYPE: Conference

LANGUAGE: English

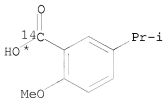
GI



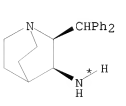
I

AB Bulk CJ-11,974-01 (I) is stable as a drug substance but degrades over time in some solid dosage formulations. Minor impurities were identified as synthetic intermediates and a major degradant has a mol. weight of M+32 by mass spectral anal., suggesting the addition of two oxygen atoms. Using solution phase hydrogen/deuterium exchange and HPLC/ESI/MS/MS techniques, the degradation product was identified as the benzyl hydroperoxide derivative The [14C]CJ-11,974-01 in ethanol solution is quite stable but is unstable as a solid, degrading to the same M+32 degradation product over a relatively short period of time. Storage of solid [14C]CJ-11,974-01 under inert atmospheric or at lower temps. did not considerably slow the degradation The carbon-14 labeled degradant was isolated by normal and reverse phase chromatogs. and identified by NMR spectroscopy and MS as the iso-Pr peroxide.

RX(2) OF 6 ...C + E ===> F...

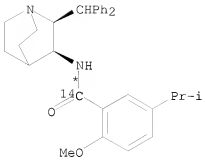


C



E

(2)



F
YIELD 88%

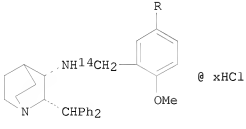
RX(2) RCT C 475146-68-0

STAGE(1)
RGT G 79-37-8 (COCl)2
SOL 75-09-2 CH2Cl2

STAGE(2)
RCT E 142035-23-2

PRO F 475146-69-1
NTE radiochem.
REFERENCE COUNT: 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

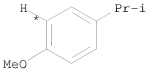
L10 ANSWER 4 OF 7 CASREACT COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 134:266175 CASREACT
TITLE: Synthesis and stability of substance P antagonists [14C]CJ-11,974-01 and [14C]CJ-11,972-01
AUTHOR(S): Zandi, K. S.; Miller, S. A.; McCarthy, K. E.; Massefski, W. W.; Kamel, A.
CORPORATE SOURCE: Radiochemical Synthesis, Pfizer Central Research, Groton, CT, 06340, USA
SOURCE: Isotope Production and Applications in the 21st Century, Proceedings of the International Conference on Isotopes, 3rd, Vancouver, BC, Canada, Sept. 6-10, 1999 (2000), Meeting Date 1999, 400-402. Editor(s): Stevenson, Nigel R. World Scientific Publishing Co. Pte. Ltd.: Singapore, Singapore.
CODEN: 69ATWE
DOCUMENT TYPE: Conference
LANGUAGE: English
GI



I

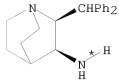
AB CJ-11,974-01 (I, R = CHMe2) and CJ-11,972-01 (I, R = CMe3) are structurally related substance P antagonists currently in development (-01 indicates the HCl salt). The synthesis of radiolabeled analogs was completed to aid in full ADME characterization. A straightforward route to both compds. was developed via directed lithiation/metal halogen exchange and carbonation. Conversion to the benzylic amine was accomplished by one of two methods. In the case of [14C]CJ-11,974-01, the carboxylic acid chloride was treated with a chiral amine followed by amide reduction, and for [14C]CJ-11,972-01, conversion to the aldehyde was followed by reductive amination. While both [14C]CJ-11,974-01 and [14C]CJ-11,972-01 are quite stable in solution, when stored as a solid, [14C]CJ-11,974-01 degrades to one major degradation product over a relatively short time period. The carbon-14 labeled degradation product was isolated from low specific activity material and identified by HPLC/MS/MS and NMR to be an iso-Pr peroxide. Studies were performed to identify the factors responsible for the oxidative degradation of [14C]CJ-11,974-01, which included salt form, storage conditions and salt formation solvent. Of all the variables studied over a three week period, only a change in the salt form prevented this oxidative degradation

RX(1) OF 2 A + B + C ===> D



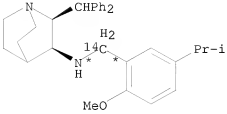
A

B



C

(1)



●x HCl

D

RX(1) RCT A 4132-48-3

STAGE(1)
RGT E 109-72-8 BuLi
SOL 109-99-9 THF

STAGE(2)
RCT B 51-90-1

STAGE(3)
RGT F 79-37-8 (COCl)2

STAGE(4)
RCT C 142035-23-2
RGT G 121-44-8 Et3N

STAGE(5)
RGT H 14044-65-6 BH3-THF

PRO D 331676-67-6

L10 ANSWER 5 OF 7 CASREACT COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 125:49301 CASREACT
TITLE: Preparation of quinuclidine derivatives as substance P antagonists
INVENTOR(S): Lowe, John Adams
PATENT ASSIGNEE(S): Pfizer Inc., India
SOURCE: Indian, 69 pp.
CODEN: INXXAP
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
IN 173570	A1	19940604	IN 1989-DE1094	19891123

PRIORITY APPLN. INFO.: IN 1989-DE1094 19891123
OTHER SOURCE(S): MARPAT 125:49301
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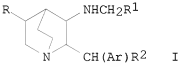
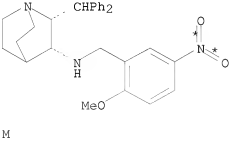


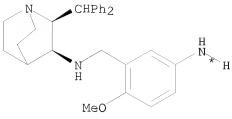
I

AB Quinuclidine derivs. [I; Ar = thienyl, Ph, halophenyl; R = H, C1-4 alkyl; R1 = C5-7 cycloalkyl, norbornyl, pyrrolyl, 2,3-dihydrobenzofuranyl, (alkoxy)thienyl, (hydroxy)pyridyl, quinolinyl, indolyl, (alkoxy)naphthyl, biphenyl, 2,3-methylenedioxyphenyl, substituted Ph, etc.; R2 = branched alkyl or alkenyl, C5-7 cycloalkyl, furyl, thienyl, (substituted) Ph, phenylalkyl, C1-3 alkoxy, etc.] are prepared for use as substance P antagonists for treatment of gastrointestinal and central nervous (psychotic) disorders, inflammatory diseases, pain, and migraine. I are prepared by reduction of the corresponding quinuclidine imine or amide. Thus, 3-keto-2-benzhydrylquinuclidine condensed with cyclohexylmethylamine to form an imine, which was reduced with 9-borabicyclononane in THF to cis-3-(cyclohexylmethylamino)-2-benzhydrylquinuclidine.

RX(3) OF 42 ...M ===> N



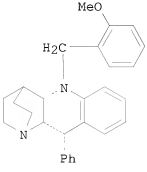
M



N
YIELD 46%

RX(3) RCT M 177746-08-6
RGT O 540-69-2 Ammonium formate
PRO N 160551-65-5
CAT 7440-05-3 Pd
SOL 64-17-5 EtOH

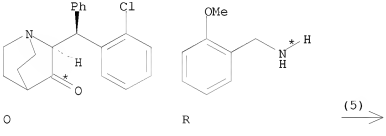
L10 ANSWER 6 OF 7 CASREACT COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 122:9904 CASREACT
TITLE: Synthesis of a benzo[b]-1,5-naphthyridine derivative as a potential constrained NK1 receptor antagonist
AUTHOR(S): Viti, Giovanni; Giannotti, Danilo; Nannicini, Rossano; Balacco, Giuseppe; Pestellini, Vittorio
CORPORATE SOURCE: Chem. Res. Dep., Firenze, 50131, Italy
SOURCE: Tetrahedron Letters (1994), 35(32), 5939-42
CODEN: TELEAY; ISSN: 0040-4039
DOCUMENT TYPE: Journal
LANGUAGE: English
GI



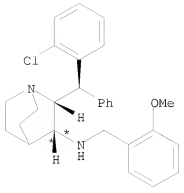
I

AB A short synthesis of a cyclic constrained analog I of the potent Substance P antagonist (±)-CP-96345 is described. The key feature is the formation of the benzo[b]-1,5-naphthyridine system at the very last step of the synthesis through an intramol. arylation of an amine promoted by a strong base. If the tricyclic system was synthesized first, 2-methoxybenzylation of both the nitrogen atoms occurred.

RX(5) OF 9 ...O + R ===> A...



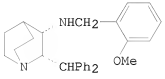
O R



A
YIELD 37%

RX(5) RCT O 159553-07-8, R 6850-57-3
RGT S 16853-85-3 LiAlH4
PRO A 159553-08-9

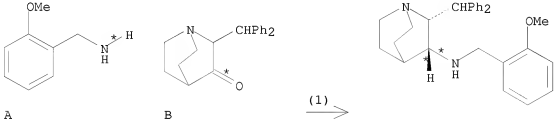
L10 ANSWER 7 OF 7 CASREACT COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 117:48289 CASREACT
TITLE: The discovery of (2S,3S)-cis-2-(diphenylmethyl)-N-[(2-methoxyphenyl)methyl]-1-azabicyclo[2.2.2]octan-3-amine as a novel, nonpeptide substance P antagonist
AUTHOR(S): Lowe, John A., III; Drozda, Susan E.; Snider, R. Michael; Longo, Kelly P.; Zorn, Stevin H.; Morrone, Jean; Jackson, Elisa R.; McLean, Stafford; Bryce, Dianne K.; et al.
CORPORATE SOURCE: Cent. Res. Div., Pfizer, Inc., Groton, CT, 06340, USA
SOURCE: Journal of Medicinal Chemistry (1992), 35(14), 2591-600
CODEN: JMCMAR; ISSN: 0022-2623
DOCUMENT TYPE: Journal
LANGUAGE: English
GI



I

AB The structure-activity relationship of a series of quinuclidines is described which culminated in the first potent, selective, nonpeptide substance P (SP) antagonist, (2S,3S)-cis-2-(diphenylmethyl)-N-[(2-methoxyphenyl)methyl]-1-azabicyclo[2.2.2]octan-3-amine, (I; CP-96,345). I is a potent displacer of [3H]SP binding in human IM-9 cells and blocks SP-induced and capsaicin-induced plasma extravasation, as well as SP-induced salivation in the rat in vivo. I may both help to further the understanding of the interactions of small mols. with peptide receptors and serve to evaluate the therapeutic potential of a SP antagonist.

RX(1) OF 35 A + B ===> C...



A B C
YIELD 71%

RX(1) RCT A 6850-57-3, B 32531-66-1
PRO C 135095-42-0
SOL 108-88-3 PhMe, 109-99-9 THF

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L2             0 S L1 FULL
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L4             0 S L3 FULL
L5               STRUCTURE UPLOADED
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L7               STRUCTURE UPLOADED
L8           335 S L7 FULL

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L10            7 S L9 FULL

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DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)          SINCE FILE      TOTAL
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